

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
 Data File : BM025817.D
 Acq On : 27 Mar 2020 01:31
 Operator : CG/JU
 Sample : L1968-13
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CODE5

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.987	341	357	366	RBV8	68460990	432535381	98.54%	25.402%
2	6.269	569	575	582	rVB	294828	454725	0.10%	0.027%
3	6.934	677	688	699	rBV	395824	872106	0.20%	0.051%
4	7.116	710	719	731	rVB	559833	1056971	0.24%	0.062%
5	7.487	770	782	793	rBV	25707199	75242239	17.14%	4.419%
6	7.686	793	816	821	RBV6	73449628	359285170	81.85%	21.100%
7	8.028	850	874	877	RBV7	73575353	438931087	100.00%	25.777%
8	8.281	911	917	928	rVB	525381	818462	0.19%	0.048%
9	8.716	985	991	994	rBV	658732	1138509	0.26%	0.067%
10	8.769	994	1000	1012	rVB	5959523	9733336	2.22%	0.572%
11	9.051	1039	1048	1059	rBV	1109594	1760339	0.40%	0.103%
12	9.357	1096	1100	1106	rVV	551382	905266	0.21%	0.053%
13	9.433	1106	1113	1120	rVV	586237	1078385	0.25%	0.063%
14	9.969	1196	1204	1215	rVB	853345	1465250	0.33%	0.086%
15	10.269	1237	1255	1259	RBV3	70102977	234653198	53.46%	13.781%
16	10.363	1265	1271	1275	rBV	990942	1427828	0.33%	0.084%
17	10.751	1323	1337	1342	rBV	38016219	73501068	16.75%	4.317%
18	10.933	1359	1368	1379	rBV	2848145	4581549	1.04%	0.269%
19	11.145	1396	1404	1411	rBV	528098	891051	0.20%	0.052%
20	12.380	1609	1614	1622	rVB	1964548	3044946	0.69%	0.179%
21	12.998	1712	1719	1727	rVB	8081102	11914103	2.71%	0.700%
22	13.621	1814	1825	1830	rBV	1731930	2586057	0.59%	0.152%
23	13.892	1866	1871	1881	rVB	2001036	2801588	0.64%	0.165%
24	14.204	1917	1924	1932	rBV	1348152	1935726	0.44%	0.114%
25	15.204	2087	2094	2100	rBV	2572406	3524488	0.80%	0.207%
26	15.327	2108	2115	2124	rBV	779637	1085067	0.25%	0.064%
27	16.951	2384	2391	2397	rBV	1702541	2296033	0.52%	0.135%
28	17.051	2401	2408	2417	rVV	2797336	3867720	0.88%	0.227%
29	18.327	2620	2625	2633	rVV	7496099	9758812	2.22%	0.573%
30	18.451	2641	2646	2654	rVB2	318888	444041	0.10%	0.026%
31	18.656	2677	2681	2685	rVV	450023	654181	0.15%	0.038%
32	19.345	2792	2798	2806	rBV	3424442	4629339	1.05%	0.272%
33	20.886	3055	3060	3066	rBV	430994	568496	0.13%	0.033%
34	21.139	3098	3103	3112	rBV	2270502	2805181	0.64%	0.165%

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Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	23.156	3439	3446	3453	rBV	3077244	5185989	1.18%	0.305%
36	23.285	3462	3468	3477	rVB2	1706837	2957052	0.67%	0.174%
37	24.080	3597	3603	3610	rBV2	276604	520934	0.12%	0.031%
38	25.579	3851	3858	3867	rVB	495045	1203402	0.27%	0.071%
39	27.603	4196	4202	4217	rVB5	207333	674195	0.15%	0.040%

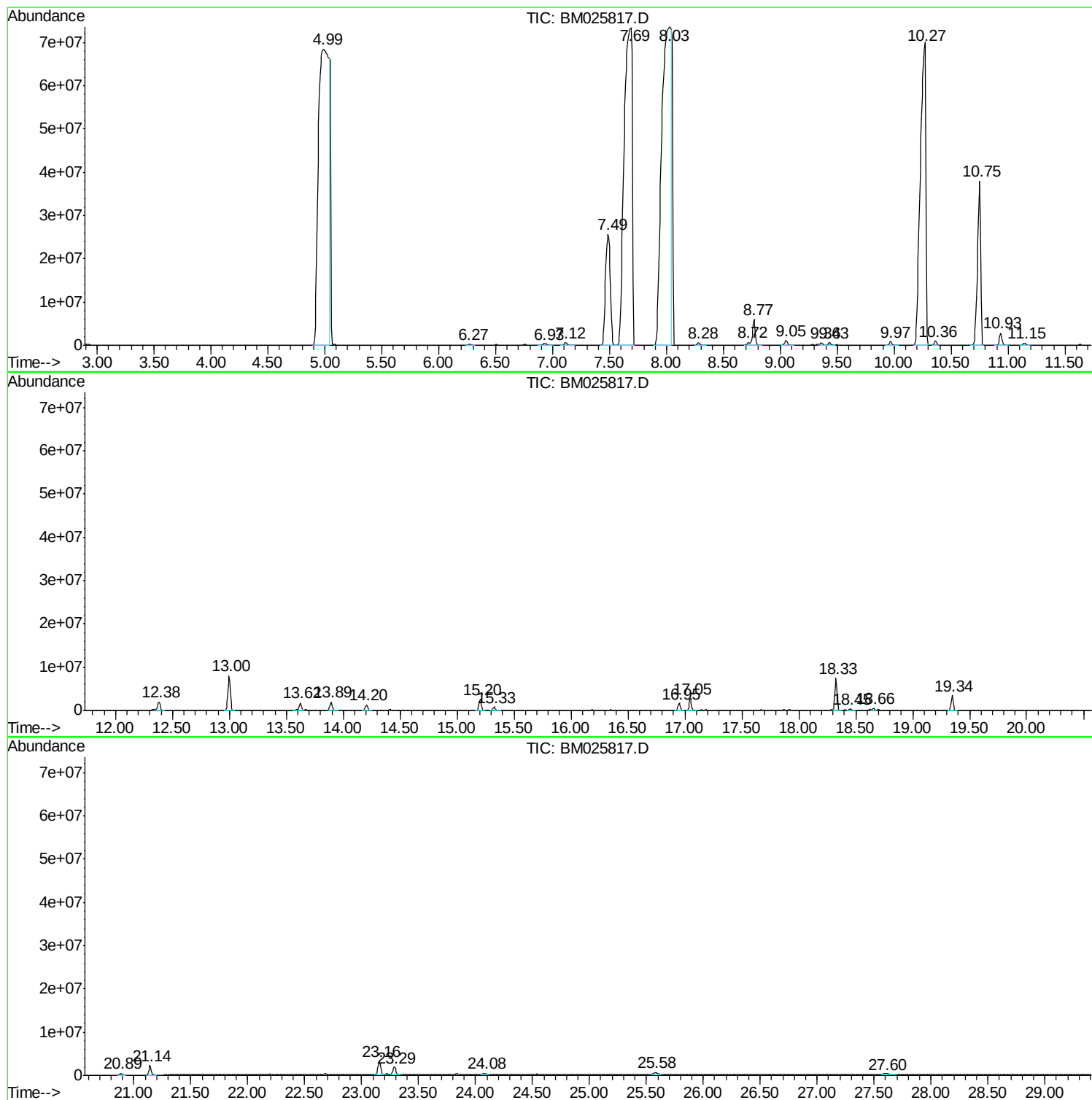
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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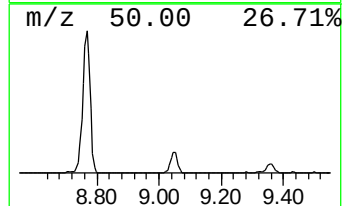
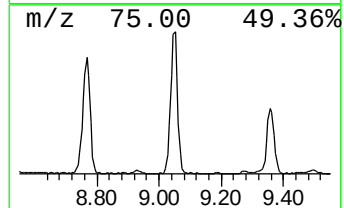
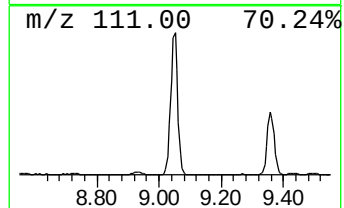
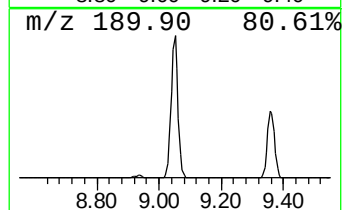
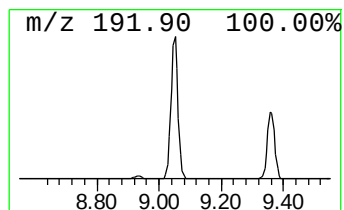
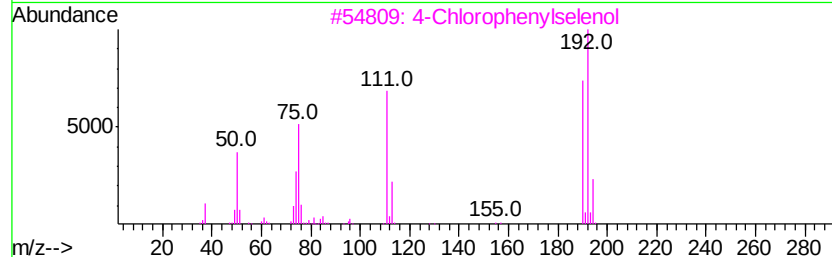
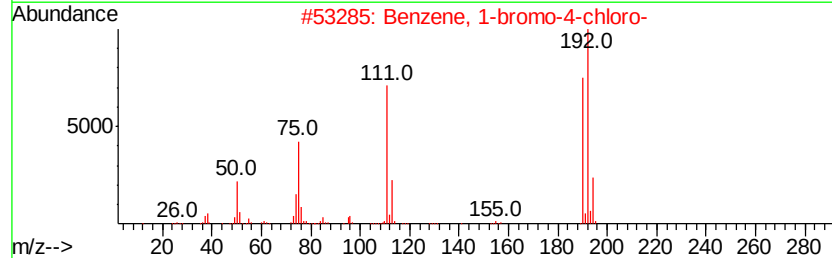
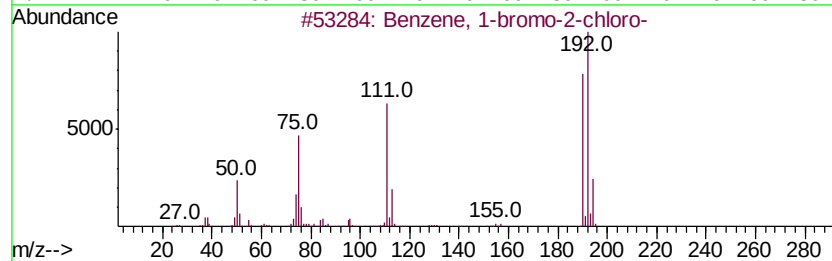
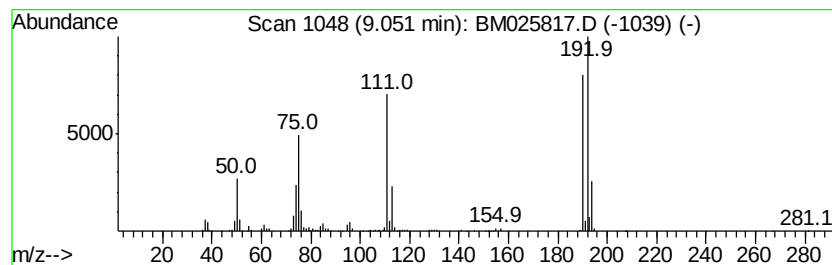
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-bromo-2-chloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.05	24.66 ng/ul	1760340	Napthalene-d8	10.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
2		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
3		4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	97
4		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96
5		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	95



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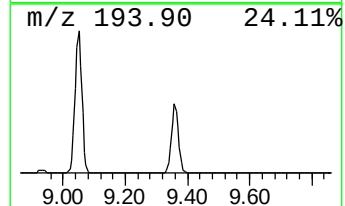
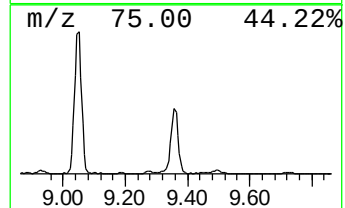
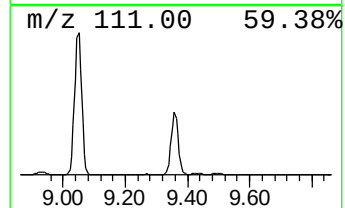
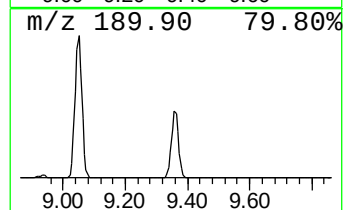
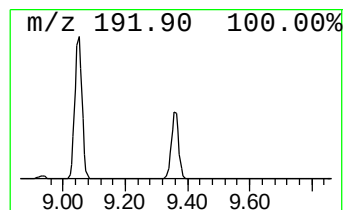
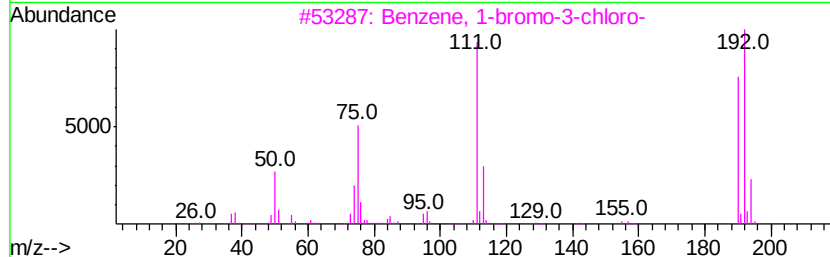
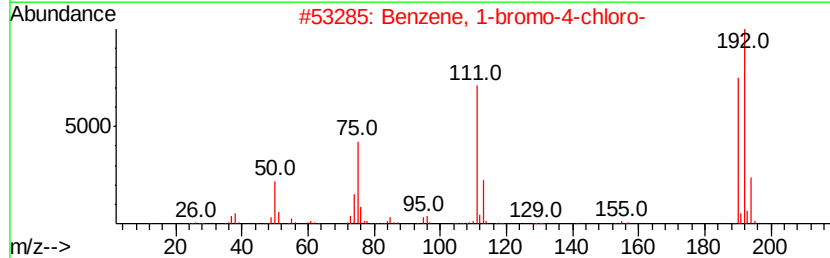
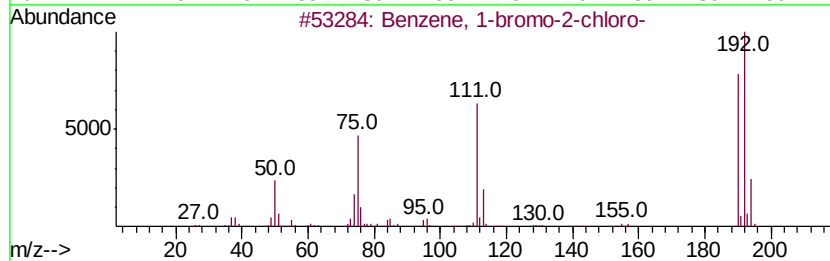
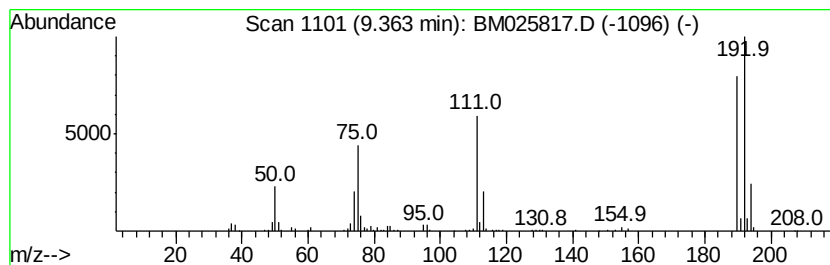
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-bromo-4-chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.36	12.68 ng/ul	905266	Napthalene-d8	10.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	99
2		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
3		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
4		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	95
5		Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	95



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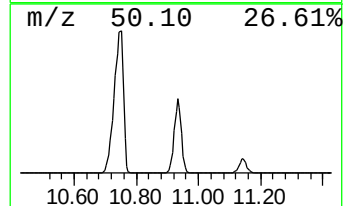
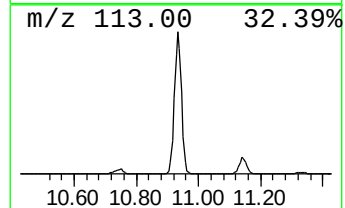
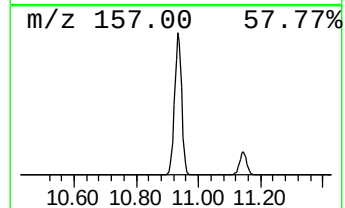
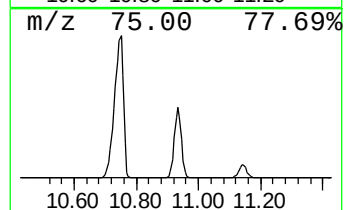
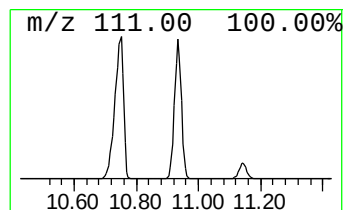
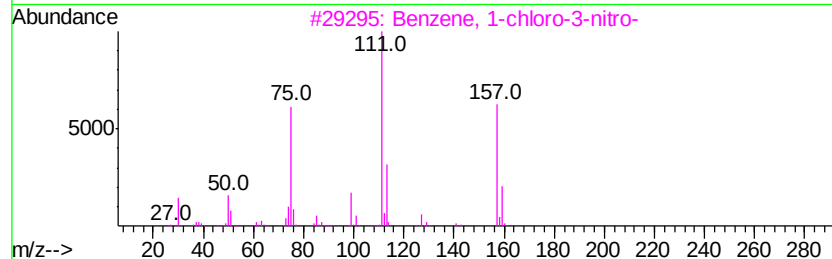
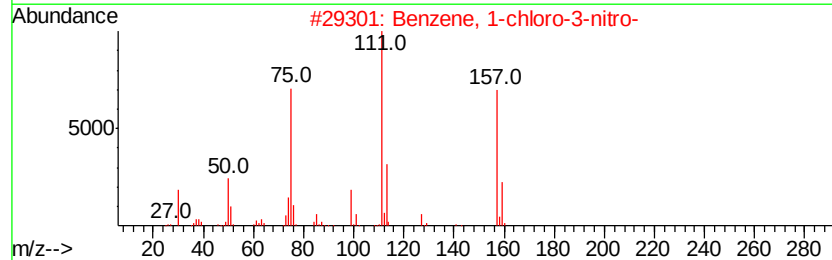
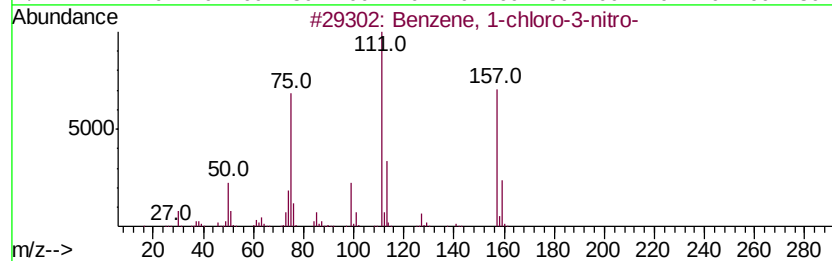
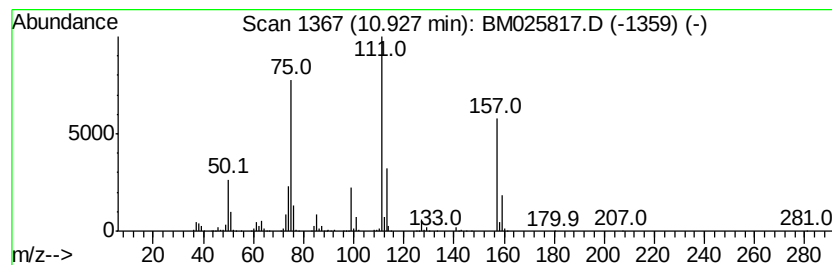
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1-chloro-3-nitro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.93	64.18 ng/ul	4581550	Napthalene-d8	10.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	98
2		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	97
3		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	96
4		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	91
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	91



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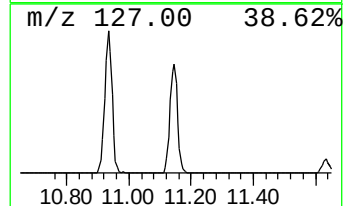
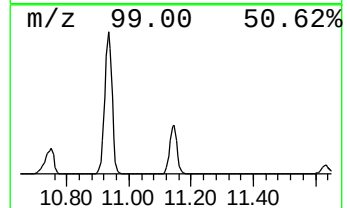
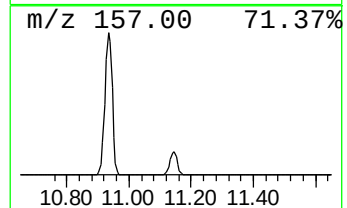
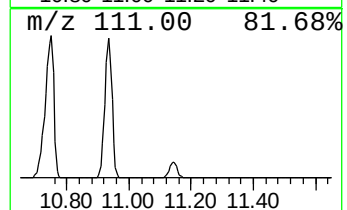
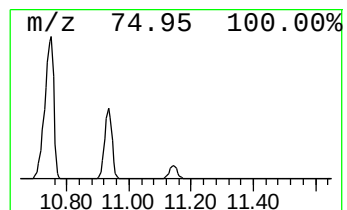
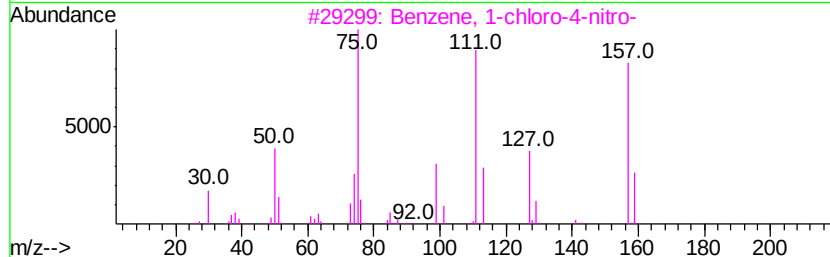
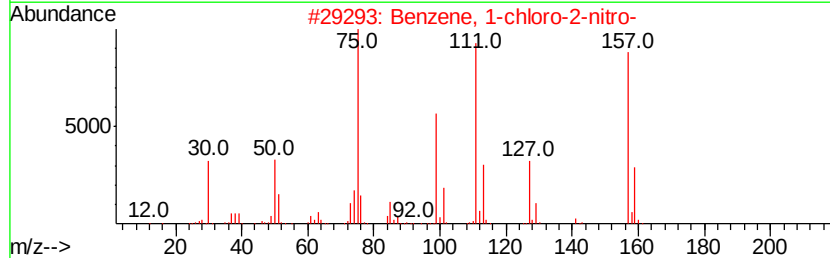
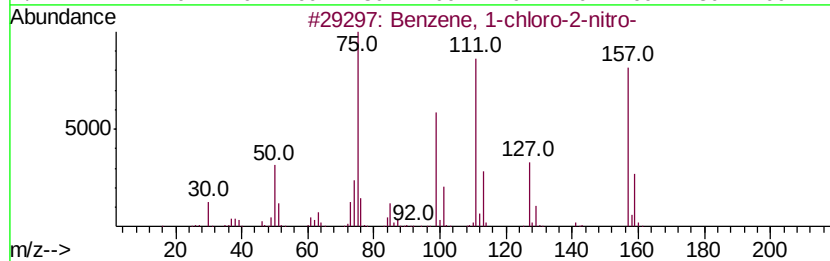
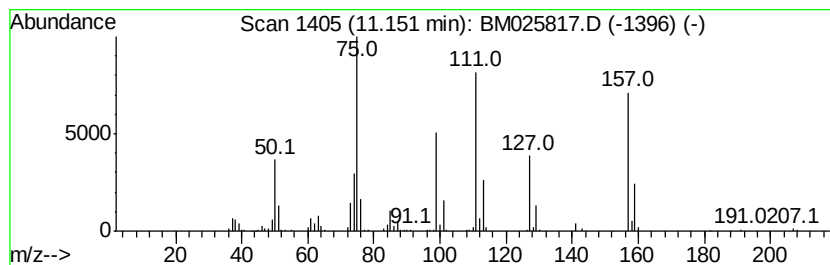
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 1-chloro-2-nitro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.15	12.48 ng/ul	891051	Napthalene-d8	10.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	99
2		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	98
3		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	98
4		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	97
5		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97



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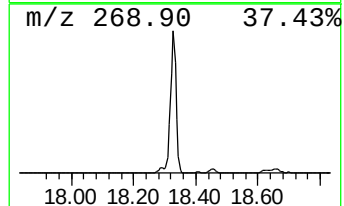
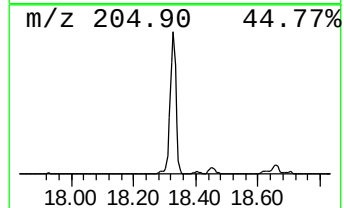
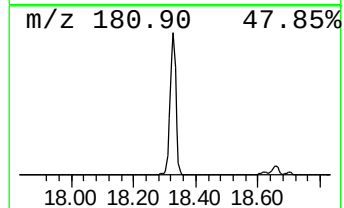
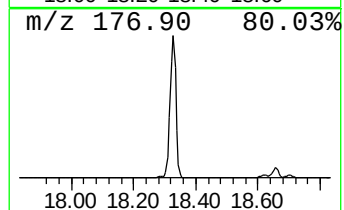
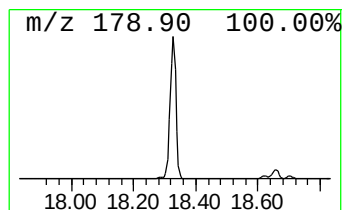
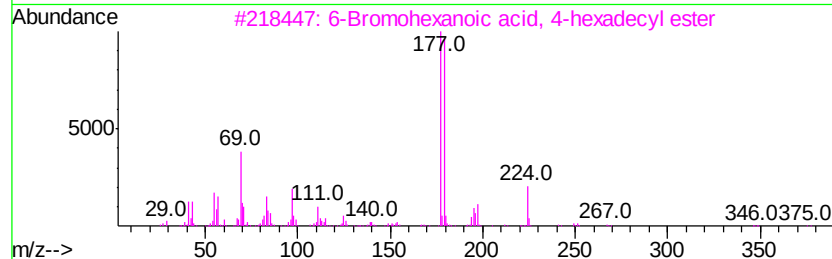
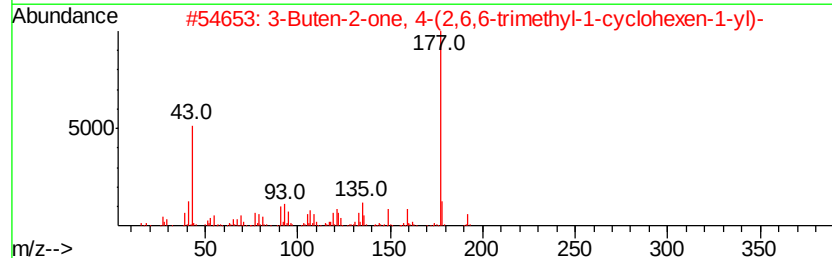
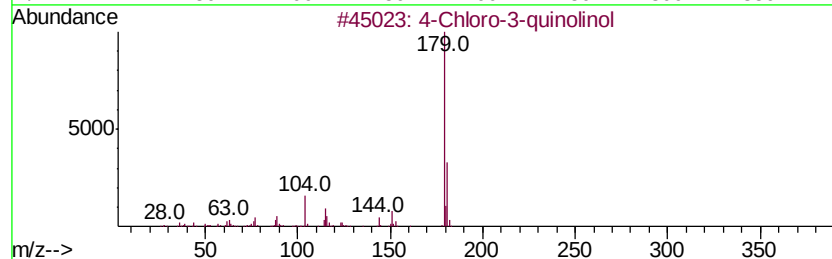
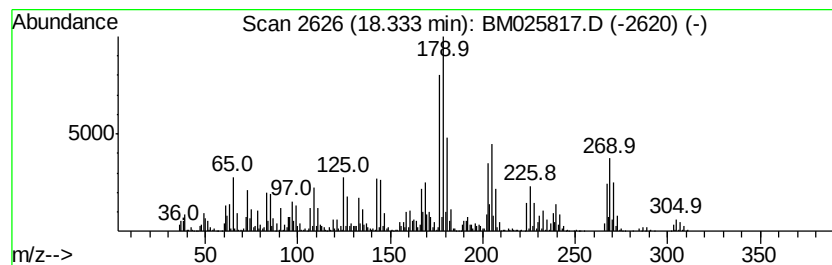
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Peak Number 11 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.33	85.01 ng/ul	9758810	Phenanthrene-d10	16.95
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		4-Chloro-3-quinolinol	179 C9H6ClNO	032435-60-2 25
2		3-Buten-2-one, 4-(2,6,6-trimethy...	192 C13H20O	014901-07-6 25
3		6-Bromohexanoic acid, 4-hexadecy...	418 C22H43BrO2	1000282-90-6 25
4		4-Butylbenzoic acid, nonyl ester	304 C20H32O2	1000292-20-6 15
5		Octamethyl-1,4-cyclohexadiene	192 C14H24	172684-93-4 15



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025817.D
Acq On : 27 Mar 2020 01:31
Operator : CG/JU
Sample : L1968-13
Misc :
ALS Vial : 17 Sample Multiplier: 1

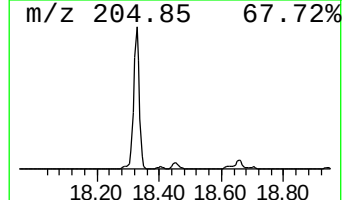
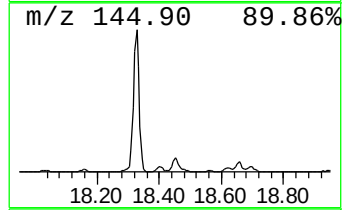
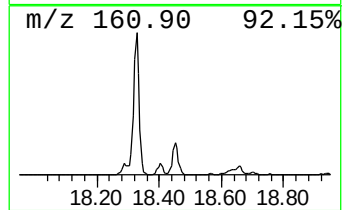
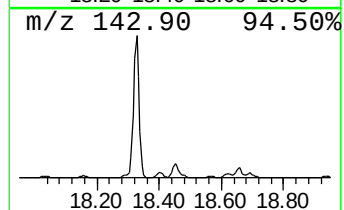
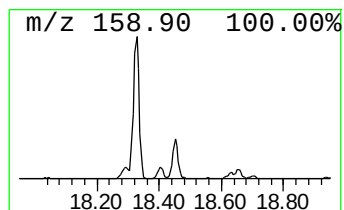
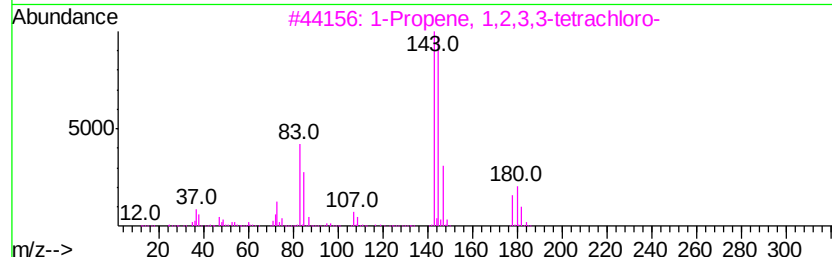
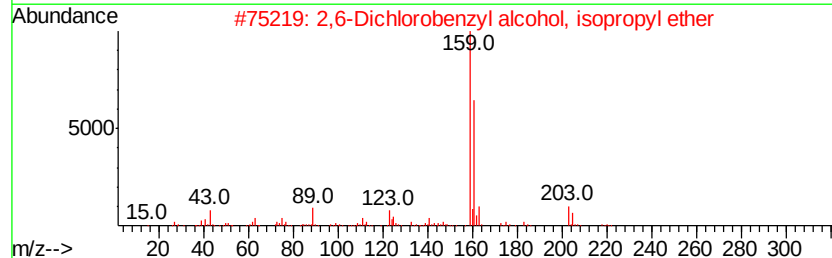
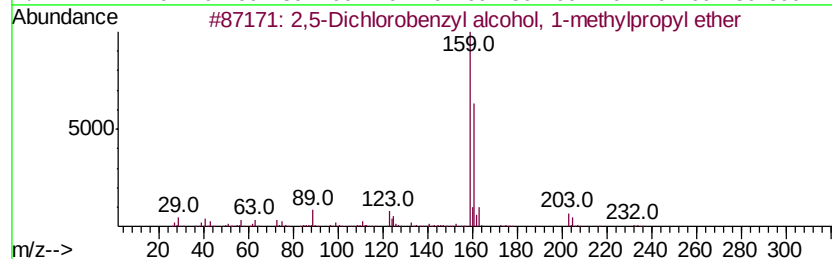
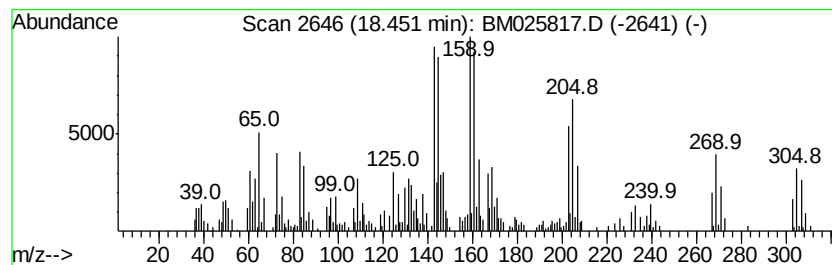
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown-02 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.45	3.87 ng/ul	444041	Phenanthrene-d10	16.95		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,5-Dichlorobenzyl alcohol, 1-me...	232	C11H14Cl2O	1000378-11-8	22	
2	2,6-Dichlorobenzyl alcohol, isop...	218	C10H12Cl2O	1000378-12-8	22	
3	1-Propene, 1,2,3,3-tetrachloro-	178	C3H2Cl4	020589-85-9	15	
4	Benzene, 2-(bromomethyl)-1,3-dic...	238	C7H5BrCl2	020443-98-5	11	
5	Acetamide, 2-(2-ethyl-1-benzimid...	203	C11H13N3O	054980-94-8	11	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
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Acq On : 27 Mar 2020 01:31
Operator : CG/JU
Sample : L1968-13
Misc :
ALS Vial : 17 Sample Multiplier: 1

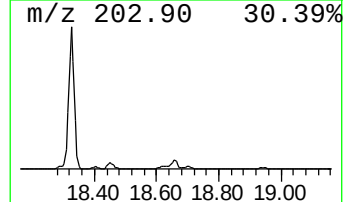
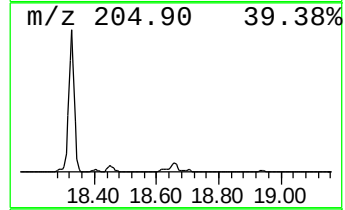
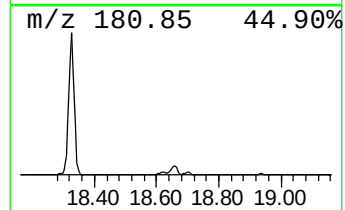
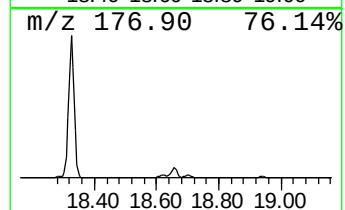
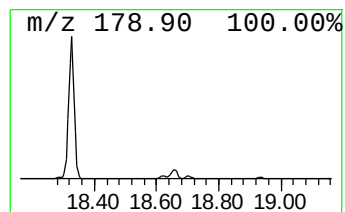
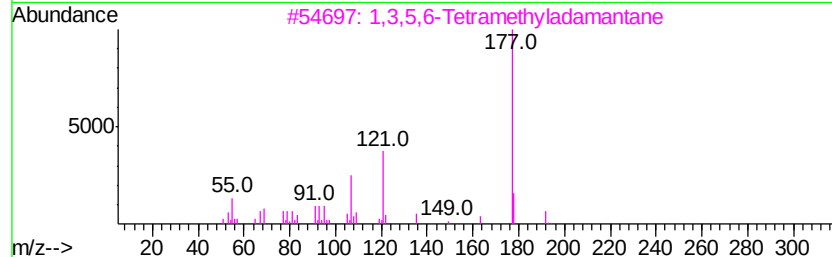
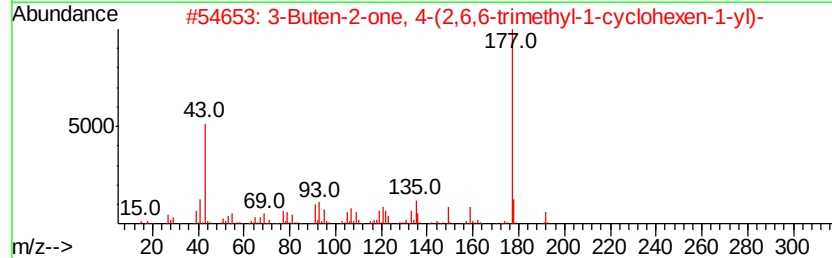
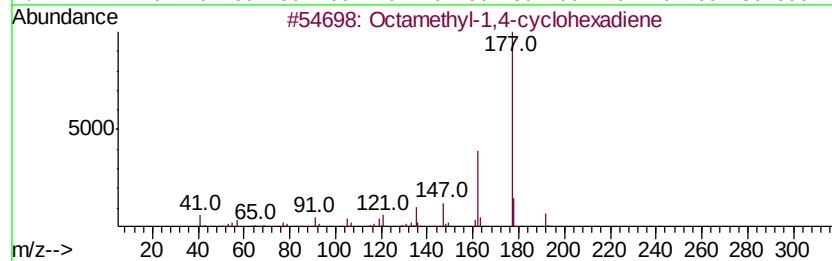
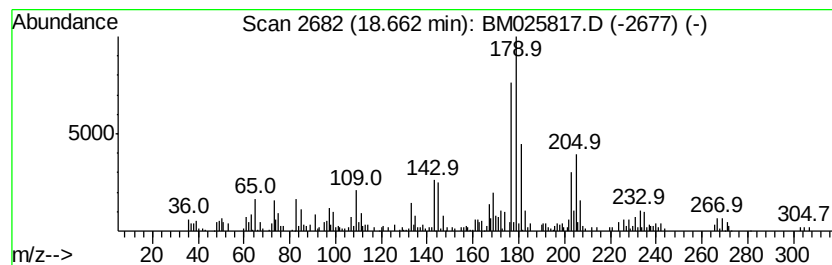
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown-03 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.66	5.70 ng/ul	654181	Phenanthrene-d10	16.95	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octamethyl-1,4-cyclohexadiene	192	C14H24	172684-93-4	25
2	3-Buten-2-one, 4-(2,6,6-trimethyl...	192	C13H20O	014901-07-6	15
3	1,3,5,6-Tetramethyladamantane	192	C14H24	034694-70-7	15
4	1,3,5,7-Tetramethyl-adamantane	192	C14H24	001687-36-1	11
5	Anthracene, 9,10-dihydro-9-(3-hy...	238	C17H18O	1000154-64-8	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025817.D
Acq On : 27 Mar 2020 01:31
Operator : CG/JU
Sample : L1968-13
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
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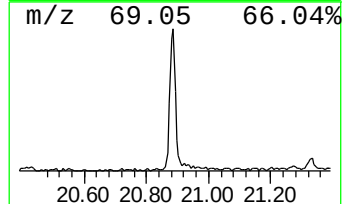
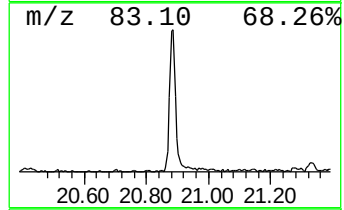
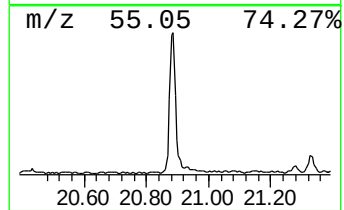
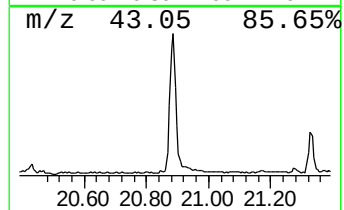
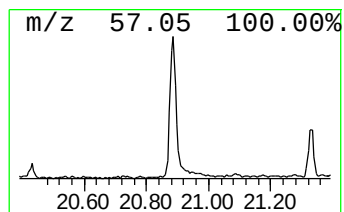
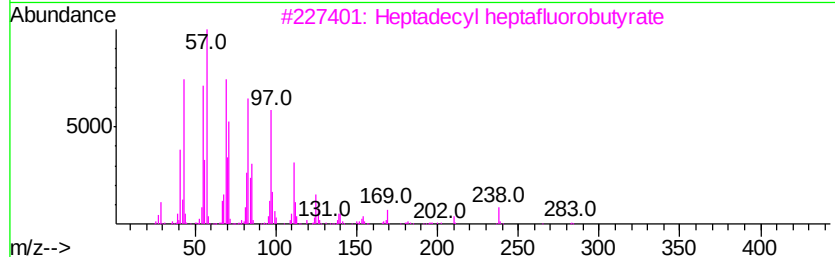
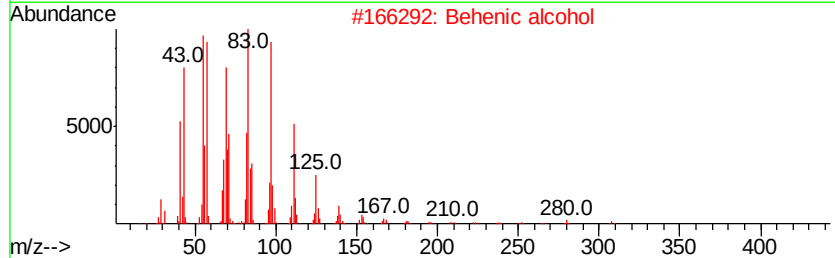
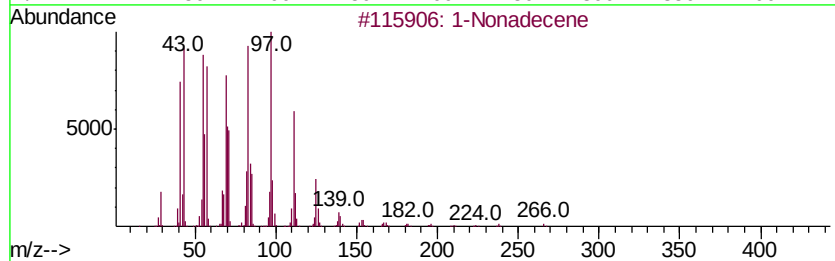
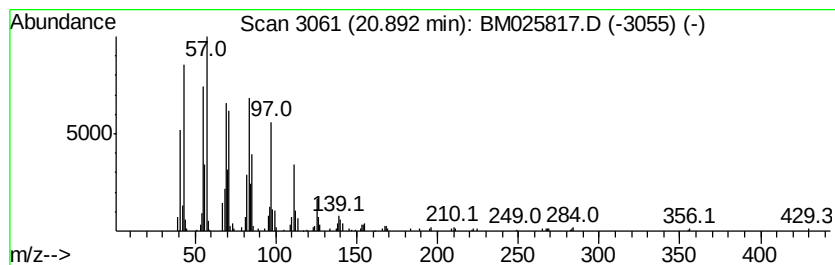
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.89	4.05 ng/ul	568496	Chrysene-d12	21.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Nonadecene	266	C19H38	018435-45-5	95
2			Behenic alcohol	326	C22H46O	000661-19-8	95
3			Heptadecyl heptafluorobutvrate	452	C21H35F7O2	959085-66-6	94
4			n-Nonadecanol-1	284	C19H40O	001454-84-8	93
5			n-Tetracosanol-1	354	C24H50O	000506-51-4	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025817.D
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Operator : CG/JU
Sample : L1968-13
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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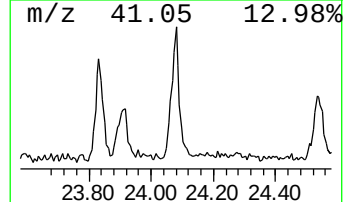
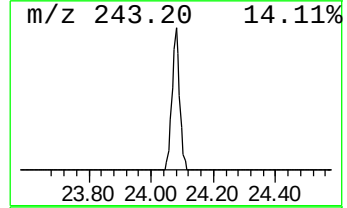
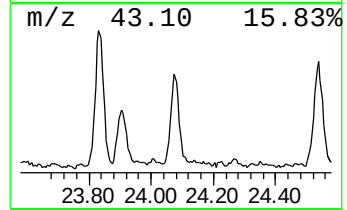
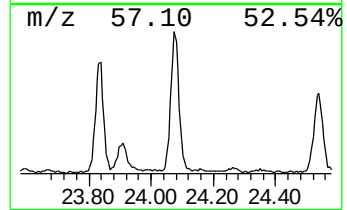
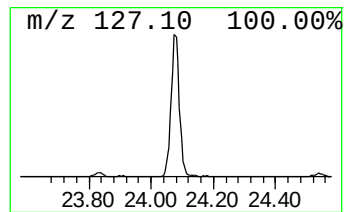
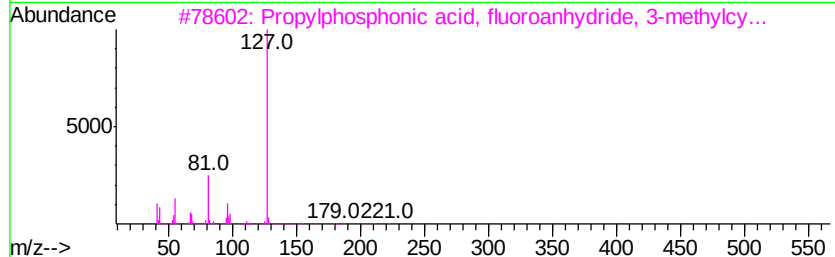
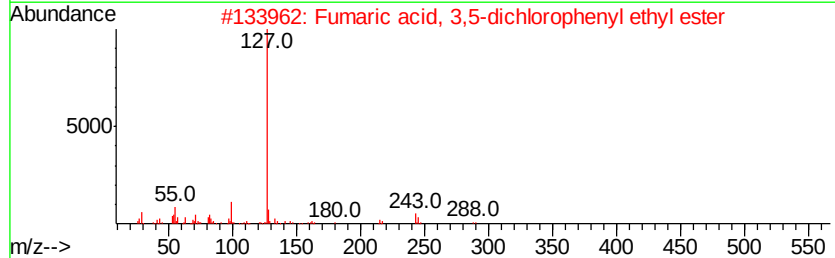
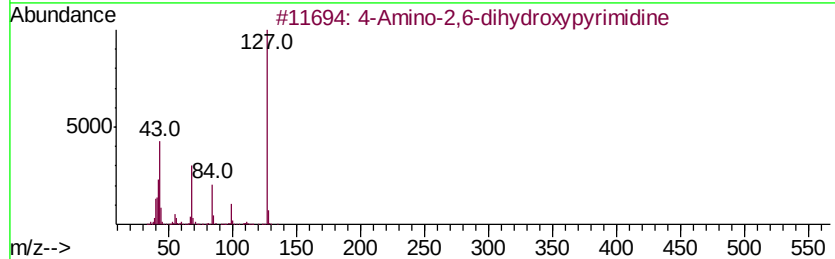
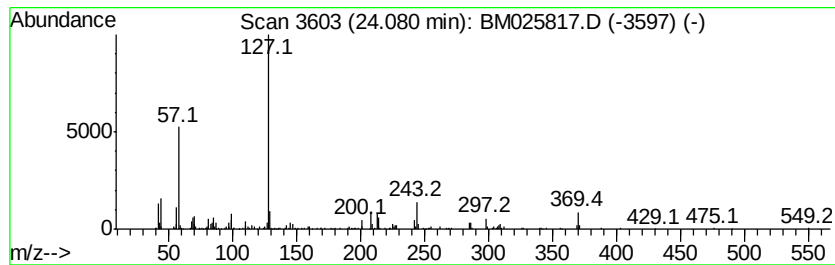
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 4-Amino-2,6-dihydroxypyrimi... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.08	3.52 ng/ul	520934	Perylene-d12	23.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Amino-2,6-dihydroxypyrimidine	127	C4H5N3O2	000873-83-6	53
2		Fumaric acid, 3,5-dichlorophenyl...	288	C12H10Cl2O4	1000348-23-1	50
3		Propylphosphonic acid, fluoroanh...	222	C10H20F02P	1000273-52-9	50
4		Phosphonofluoridic acid, (1-meth...	222	C10H20F02P	333416-73-2	50
5		(Chlorophenoxymethylenimino)sulf...	281	C7H5ClF5NOS	081439-21-6	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025817.D
Acq On : 27 Mar 2020 01:31
Operator : CG/JU
Sample : L1968-13
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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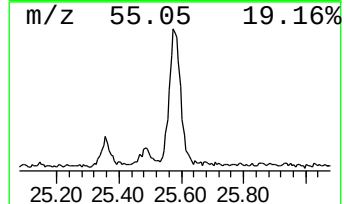
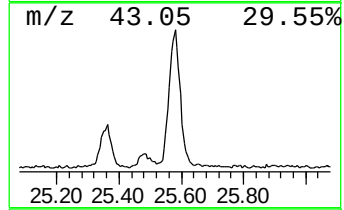
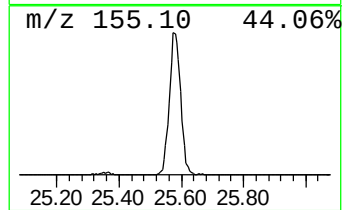
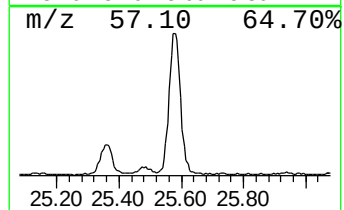
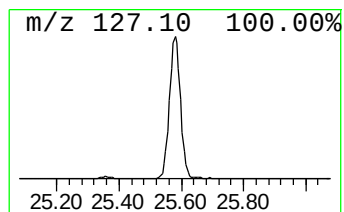
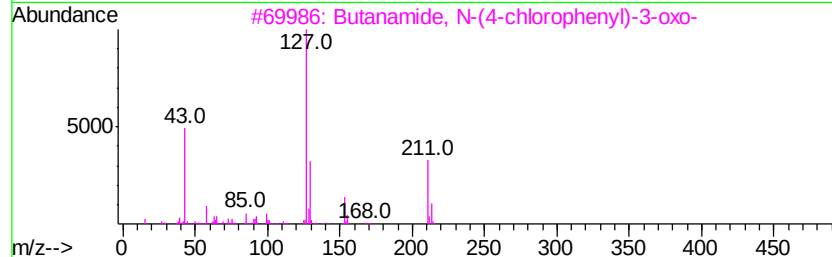
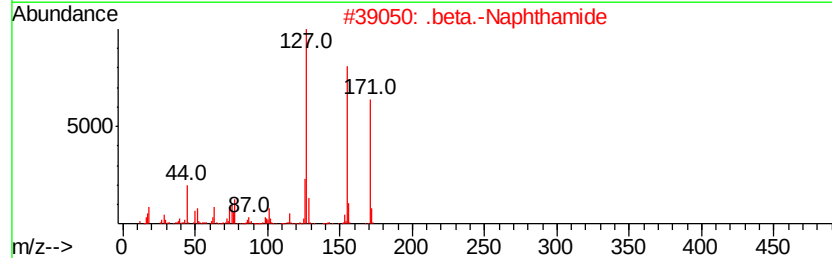
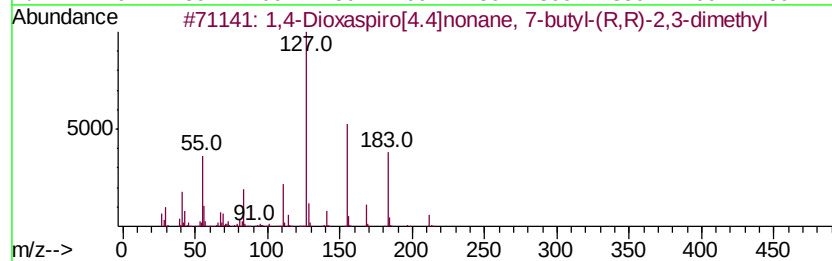
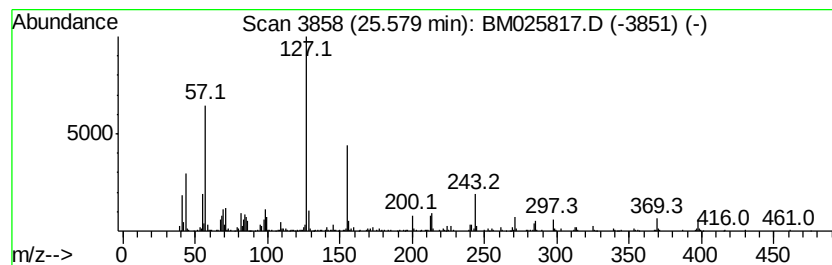
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 1,4-Dioxaspiro[4.4]nonane, ... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.58	8.14 ng/ul	1203400	Perylene-d12	23.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Dioxaspiro[4.4]nonane, 7-but...	212	C13H24O2	104807-77-4	64
2			.beta.-Naphthamide	171	C11H9NO	002243-82-5	59
3			Butanamide, N-(4-chlorophenyl)-3...	211	C10H10ClNO2	000101-92-8	43
4			2-Hydroxy-4-hydroxylaminopyrimidine	127	C4H5N3O2	020555-88-8	43
5			2,2'-Dithiodi[3-pyridinol]	252	C10H8N2O2S2	068750-09-4	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025817.D
Acq On : 27 Mar 2020 01:31
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Sample : L1968-13
Misc :
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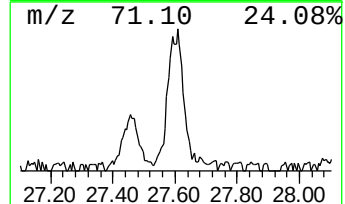
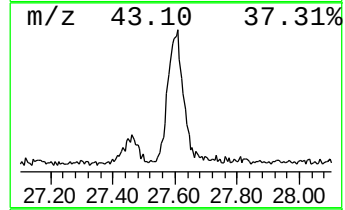
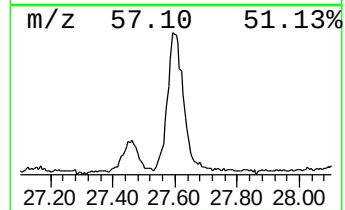
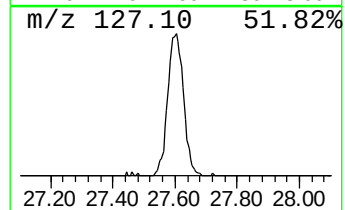
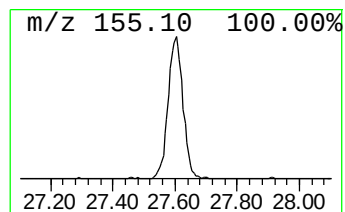
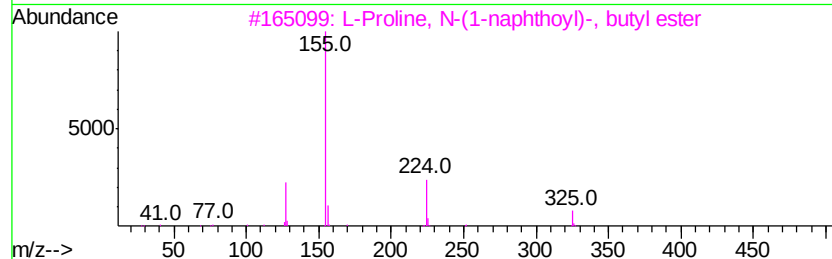
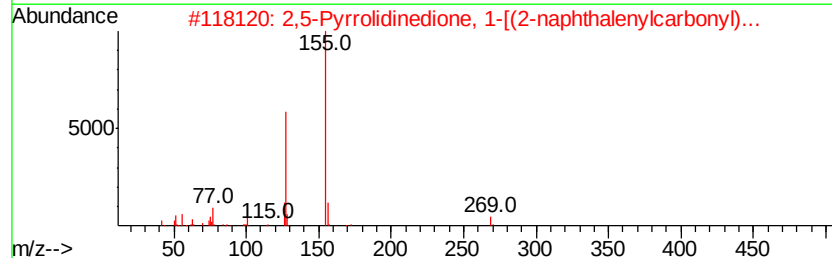
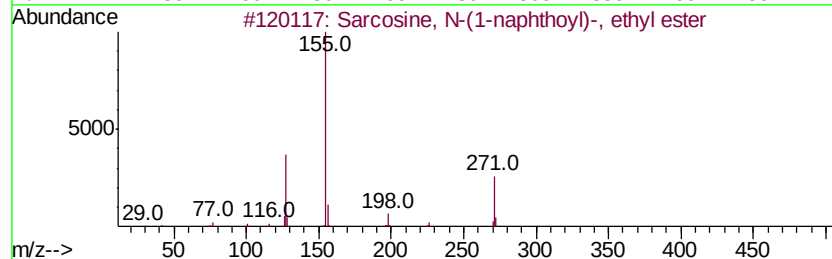
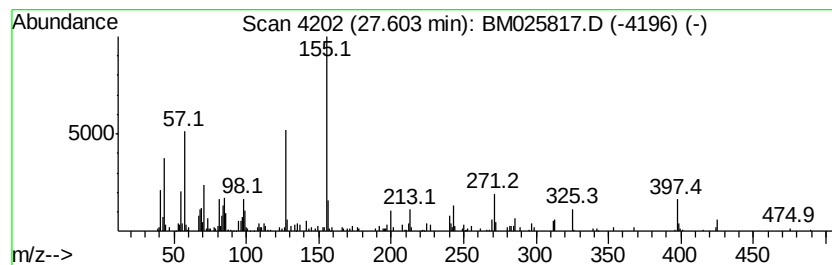
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Sarcosine, N-(1-naphthoyl)-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.60	4.56 ng/ul	674195	Perylene-d12	23.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sarcosine, N-(1-naphthoyl)-, eth...	271	C16H17N03	1000321-39-8	53
2			2,5-Pyrrolidinedione, 1-[(2-naph...	269	C15H11N04	056374-47-1	53
3			L-Proline, N-(1-naphthoyl)-, but...	325	C20H23N03	1000346-08-5	52
4			5-Oxazolidinone, 2-(1,1-dimethyl...	297	C18H19N03	119215-56-4	49
5			L-Proline, N-(1-naphthoyl)-, iso...	325	C20H23N03	1000346-08-4	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032620\
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 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-bromo-...	9.05	24.7	ng/ul	1760340	2	10.36	1427830	20.0
Benzene, 1-bromo-...	9.36	12.7	ng/ul	905266	2	10.36	1427830	20.0
Benzene, 1-chloro...	10.93	64.2	ng/ul	4581550	2	10.36	1427830	20.0
Benzene, 1-chloro...	11.15	12.5	ng/ul	891051	2	10.36	1427830	20.0
unknown-01	18.33	85.0	ng/ul	9758810	4	16.95	2296030	20.0
unknown-02	18.45	3.9	ng/ul	444041	4	16.95	2296030	20.0
unknown-03	18.66	5.7	ng/ul	654181	4	16.95	2296030	20.0
(DEL) Alkane: Str...	20.89	4.0	ng/ul	568496	5	21.14	2805180	20.0
4-Amino-2,6-dihyd...	24.08	3.5	ng/ul	520934	6	23.29	2957050	20.0
1,4-Dioxaspiro[4...	25.58	8.1	ng/ul	1203400	6	23.29	2957050	20.0
Sarcosine, N-(1-n...	27.60	4.6	ng/ul	674195	6	23.29	2957050	20.0